R_1 is haloalkyl, hydroxyalkyl, haloalkenyl, heterocycloalkyl, unsubstituted phenyl, substituted or unsubstituted phenylamino, unsubstituted benzyl, alkoxyalkyl, poly(alkoxy)alkyl, hydroxyalkoxyalkyl, hydroxypoly(alkoxy)alkyl, haloalkoxyalkyl, halopoly(alkoxy)alkyl, or aminoalkyl.

REMARKS

I. Amendments to Specification

The title has been amended to be more descriptive of the invention to which the claims are directed.

Tables 2 and 3 have been amended to correct an inadvertent typographical error. "R" in the Tables should be "R_1". Support for these amendments is found at page 13, line 12, and at page 14, line 2.

II. Amendment to the Claims

Claim 1 has been amended herein.

III. Conclusion

Should the Examiner have any questions, comments or suggestions that would expedite the prosecution of the present case to allowance, Applicants' representative, Dr. Rose Ann Dabek, earnestly requests a telephone conference at (513) 627-8824.

Respectfully submitted,

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Date: July 18, 2001

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ATTACHMENT A
VERSION WITH MARKINGS TO SHOW CHANGES MADE

In the Specification:

At page 13, Table 2 is amended, as indicated below:

Table 2

Cpd. No.	[R] R_1	R_2	LogP
2-1	$-CH_2CH_2CH_2N(CH_2CH_3)_2$	H	1.098
2-2	$-CH_2CH_2$-morpholino	H	0.018
2-3	$-CH(CH_3)CH_2CH_3$	H	1.606
2-4	$-CH_2$-(2-tetrahydrofuryl)	H	0.613
2-5	$-CH_2CH_2CH(CH_3)_2$	H	1.920
2-6	$-CH(CH_3)CH_2CH(CH_3)_2$	H	2.333
2-7	$-CH_2CH_2C(CH_3)_3$	H	2.353
2-8	$-CH_2CH(CH_3)CH_2CH_3$	H	1.992
2-9	$-CH\ (CH_2CH_3)_2$	H	2.075
2-10	$-CH_2CH_2CH_2N(CH_3)_2$	H	0.413
2-11	$-CH_2CH_2OCH_3$	H	0.217
2-12	-NH-Ph	H	1.737
2-13	-Ph(2-OH)	H	1.779
2-14	$-CH_2CH_2N(CH_3)_2$	H	0.361
2-15	$-Ph(3-OCH_3-4-OCH_3-5-OCH_3)$	H	1.305
2-16	cyclohexyl	CH_3	2.213
2-17	$-CH_2CH_2CH_2CH_3$	CH_3	1.836
2-18	$-CH_2CH_2CH_3$	$CH_2CH_2CH_3$	2.250

At page 14, Table 3 is amended, as indicated below:

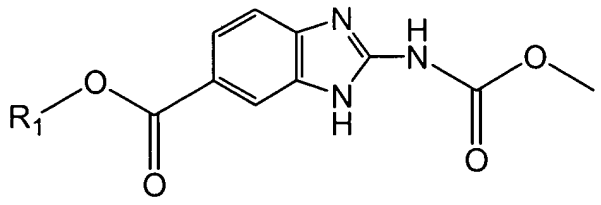
Table 3

Cpd. No.	[R] R_1	Log P
3-1	$-CH_2CH_2CH_2CH_2Cl$	2.239
3-2	$-CH_2CH_2OCH_2CH_2Cl$	1.571
3-3	$-CH_2CH=CH_2$	1.772
3-4	$-(CH_2CH_2O)_2CH_2CH_3$	1.045
3-5	$-CH_2CH_2OCH_2CH_2OH$	0.424
3-6	$-CH_2CH_2CH=CH_2$	2.024
3-7	$-CH_2Ph$	2.808
3-8	$-CH_2CH_2N(CH_3)_2$	1.011
3-9	$-CH_2CH_2CH_2Cl$	1.788
3-10	$-CH_2CH=CHCH_2OH$	1.121
3-11	$-CH_2CH_2CH_2CH_2CH_2OH$	1.488
3-12	$-CH(CH_2Cl)_2$	2.510
3-13	$-CH_2CH(CH_3)CH_2C(CH_3)_3$	3.802
3-14	$-CH_2CF_2CF_3$	2.841
3-15	$-CH(CH_2F)_2$	1.423
3-16	$-CH(CH_3)$(cyclopropyl)	2.155
3-17	$-CH_2CH_2F$	0.542
3-18	$-CH(CH_2Br)_2$	2.636
3-19	$-CH_2CH(CH_3)CH_2CH_3$	2.256
3-20	$-CH_2CH_2CH(CH_3)CH_2C(CH_3)_3$	4.126
3-21	$-CH_2CH_2CH(CH_3)CH_2CH_2CH=C(CH_3)_2$	4.048

In the Claims:

The following claim 1 is amended, as indicated below:

1. (1st Time Amended) A compound of the following formula **A-3**:



A-3

wherein,

R_1 is haloalkyl, hydroxyalkyl, haloalkenyl, heterocycloalkyl, [substituted or] unsubstituted phenyl, substituted or unsubstituted phenylamino, unsubstituted benzyl, alkoxyalkyl, poly(alkoxy)alkyl, hydroxyalkoxyalkyl, hydroxypoly(alkoxy)alkyl, haloalkoxyalkyl, halopoly(alkoxy)alkyl, or aminoalkyl.